

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
 Data File : BF133739.D
 Acq On : 13 Jun 2023 19:35
 Operator : CG\JU
 Sample : 03138-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-1

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/14/2023
 Supervised By :mohammad ahmed 06/16/2023

Quant Time: Jun 13 21:50:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 13 11:26:55 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	83915	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	264923	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	128391	20.000	ng	-0.01
64) Phenanthrene-d10	11.375	188	215944	20.000	ng	# 0.00
76) Chrysene-d12	14.039	240	156486	20.000	ng	# 0.02
86) Perylene-d12	15.492	264	133108	20.000	ng	# 0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	345413	71.659	ng	0.00
7) Phenol-d6	6.457	99	364956	58.167	ng	-0.01
23) Nitrobenzene-d5	7.393	82	225532	46.368	ng	-0.01
42) 2,4,6-Tribromophenol	10.663	330	101541	62.479	ng	0.00
45) 2-Fluorobiphenyl	9.187	172	417965	46.276	ng	-0.01
79) Terphenyl-d14	12.975	244	490246	48.476	ng	0.02
Target Compounds						Qvalue
31) Naphthalene	8.134	128	292469	22.660	ng	100
37) 2-Methylnaphthalene	8.828	142	100491	11.206	ng	98
38) 1-Methylnaphthalene	8.928	142	83448	10.112	ng	98
46) 1,1'-Biphenyl	9.292	154	31315	3.024	ng	100
49) Acenaphthylene	9.734	152	40793	3.155	ng	# 85
52) Acenaphthene	9.904	154	283752	37.497	ng	99
55) Dibenzofuran	10.075	168	221877	19.215	ng	96
58) Fluorene	10.422	166	225792	25.570	ng	98
71) Phenanthrene	11.416	178	3542373	323.330	ng	99
72) Anthracene	11.457	178	738880	64.680	ng	99
73) Carbazole	11.604	167	346693	32.585	ng	# 94
75) Fluoranthene	12.633	202	6036724	518.867	ng	97
78) Pyrene	12.869	202	5273594	361.769	ng	98
81) Benzo(a)anthracene	14.033	228	2788715	257.637	ng	97
83) Chrysene	14.074	228	2414725m	235.865	ng	
84) Bis(2-ethylhexyl)phtha...	13.998	149	32812	4.330	ng	# 93
87) Indeno(1,2,3-cd)pyrene	16.968	276	768612	83.932	ng	97
88) Benzo(b)fluoranthene	15.080	252	2260583m	279.576	ng	
89) Benzo(k)fluoranthene	15.104	252	728432m	92.482	ng	
90) Benzo(a)pyrene	15.445	252	1219837	162.554	ng	97
91) Dibenzo(a,h)anthracene	16.974	278	199436	26.306	ng	95
92) Benzo(g,h,i)perylene	17.415	276	721657	96.006	ng	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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