

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071723\
 Data File : BF134318.D
 Acq On : 17 Jul 2023 16:25
 Operator : CG\JU
 Sample : 03564-02DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-225DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/19/2023
 Supervised By :mohammad ahmed 07/19/2023

Quant Time: Jul 17 16:58:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 14 22:51:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	81421	20.000	ng	-0.01
21) Naphthalene-d8	8.116	136	307869	20.000	ng	0.00
39) Acenaphthene-d10	9.875	164	146667	20.000	ng	0.00
64) Phenanthrene-d10	11.369	188	253983	20.000	ng	0.00
76) Chrysene-d12	14.033	240	162529	20.000	ng	0.00
86) Perylene-d12	15.533	264	123998	20.000	ng	#-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	240301	49.369	ng	0.00
7) Phenol-d6	6.475	99	284189	47.476	ng	-0.01
23) Nitrobenzene-d5	7.398	82	185607	32.498	ng	-0.01
42) 2,4,6-Tribromophenol	10.669	330	81362	44.245	ng	-0.01
45) 2-Fluorobiphenyl	9.192	172	344916	34.191	ng	-0.01
79) Terphenyl-d14	12.963	244	336532	33.325	ng	-0.01
Target Compounds						Qvalue
52) Acenaphthene	9.904	154	24834	2.911	ng	99
58) Fluorene	10.422	166	42842	4.130	ng	100
71) Phenanthrene	11.392	178	368302	28.805	ng	98
72) Anthracene	11.445	178	89813	6.753	ng	98
73) Carbazole	11.604	167	36848	3.027	ng	100
75) Fluoranthene	12.592	202	615412	44.522	ng	98
78) Pyrene	12.822	202	505768	33.438	ng	100
81) Benzo(a)anthracene	14.022	228	316772	28.103	ng	99
83) Chrysene	14.057	228	265876	24.354	ng	97
87) Indeno(1,2,3-cd)pyrene	17.074	276	116343	13.522	ng	# 94
88) Benzo(b)fluoranthene	15.092	252	384223m	52.697	ng	
89) Benzo(k)fluoranthene	15.121	252	125403m	16.893	ng	
90) Benzo(a)pyrene	15.468	252	194143	27.536	ng	98
91) Dibenzo(a,h)anthracene	17.086	278	39332	5.597	ng	97
92) Benzo(g,h,i)perylene	17.545	276	106387	14.679	ng	99

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

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