

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
 Data File : BF137493.D
 Acq On : 04 Mar 2024 21:17
 Operator : CG\JU
 Sample : P1666-01 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-030124

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 03/05/2024
 Supervised By :mohammad ahmed 03/05/2024

Quant Time: Mar 04 23:20:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 04 00:26:53 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	194710	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	744062	20.000	ng	0.00
39) Acenaphthene-d10	9.886	164	374495	20.000	ng	0.00
64) Phenanthrene-d10	11.381	188	509154	20.000	ng	0.00
76) Chrysene-d12	14.045	240	365154	20.000	ng	0.00
86) Perylene-d12	15.574	264	289666	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	583520	45.389	ng	-0.01
7) Phenol-d6	6.493	99	822948	44.346	ng	-0.01
23) Nitrobenzene-d5	7.410	82	453933	28.347	ng	-0.02
42) 2,4,6-Tribromophenol	10.681	330	98926	30.080	ng	-0.01
45) 2-Fluorobiphenyl	9.204	172	854035	35.698	ng	-0.01
79) Terphenyl-d14	12.980	244	656676	24.729	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.151	128	240922	6.035	ng	98
37) 2-Methylnaphthalene	8.845	142	147349	5.790	ng	95
38) 1-Methylnaphthalene	8.939	142	135621	5.659	ng	98
52) Acenaphthene	9.916	154	177642	8.897	ng	99
55) Dibenzofuran	10.092	168	206187	6.780	ng	99
58) Fluorene	10.439	166	290233	12.427	ng	97
71) Phenanthrene	11.404	178	1173664	42.356	ng	99
72) Anthracene	11.457	178	293329	10.307	ng	98
73) Carbazole	11.616	167	238637	9.789	ng	98
75) Fluoranthene	12.604	202	512142	18.909	ng	97
78) Pyrene	12.833	202	430312	12.010	ng	98
81) Benzo(a)anthracene	14.033	228	148589	5.806	ng	96
83) Chrysene	14.069	228	136587m	5.280	ng	
88) Benzo(b)fluoranthene	15.121	252	110184m	6.430	ng	
89) Benzo(k)fluoranthene	15.145	252	42144m	2.290	ng	
90) Benzo(a)pyrene	15.504	252	76414	4.525	ng	# 82
92) Benzo(g,h,i)perylene	17.633	276	37956	2.390	ng	# 91

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

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