

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
 Data File : BP013630.D
 Acq On : 27 Jan 2023 22:16
 Operator : CG/JU
 Sample : 01221-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-10-0118

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/30/2023
 Supervised By : Jagrut Upadhyay 01/30/2023

Quant Time: Jan 29 23:55:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 27 03:27:54 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.152	152	190963	20.000	ng	0.00
21) Naphthalene-d8	10.981	136	711597	20.000	ng	0.00
39) Acenaphthene-d10	14.787	164	434425	20.000	ng	0.00
64) Phenanthrene-d10	17.545	188	950176	20.000	ng	# 0.00
76) Chrysene-d12	21.627	240	1053888	20.000	ng	0.00
86) Perylene-d12	24.215	264	1226184	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.687	112	1057965	99.278	ng	0.00
7) Phenol-d6	7.299	99	1034783	71.881	ng	0.00
23) Nitrobenzene-d5	9.328	82	935217	72.215	ng	0.00
42) 2,4,6-Tribromophenol	16.281	330	791117	119.611	ng	0.00
45) 2-Fluorobiphenyl	13.416	172	2513127	82.622	ng	0.00
79) Terphenyl-d14	20.080	244	4108593	68.795	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.322	94	235839	15.639	ng	96
17) 2-Methylphenol	8.593	107	38604	3.548	ng	96
20) 3+4-Methylphenols	8.922	107	53001	3.554	ng	92
27) 2,4-Dimethylphenol	10.169	122	80186m	7.440	ng	
31) Naphthalene	11.034	128	1943832	53.757	ng	99
36) 4-Chloro-3-methylphenol	12.240	107	178113	13.852	ng	99
38) 1-Methylnaphthalene	12.845	142	214506	8.293	ng	95
84) Bis(2-ethylhexyl)phtha...	21.516	149	18858	4.536	ng	# 83

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

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