

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051925\
 Data File : BP024696.D
 Acq On : 19 May 2025 19:40
 Operator : RC/JU
 Sample : Q2067-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 303-PPR-FRAC

Quant Time: May 20 03:46:44 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 13 16:21:39 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.657	152	135379	20.000	ng	0.00
21) Naphthalene-d8	10.428	136	558037	20.000	ng	0.00
39) Acenaphthene-d10	14.298	164	375594	20.000	ng	0.00
64) Phenanthrene-d10	17.098	188	788156	20.000	ng	0.00
76) Chrysene-d12	21.533	240	1047049	20.000	ng	0.01
86) Perylene-d12	24.804	264	1155824	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	495403	62.589	ng	0.00
7) Phenol-d6	6.869	99	382330	37.847	ng	0.02
23) Nitrobenzene-d5	8.804	82	893899	80.936	ng	0.00
42) 2,4,6-Tribromophenol	15.822	330	932116	146.389	ng	0.01
45) 2-Fluorobiphenyl	12.904	172	1930557	67.341	ng	0.00
79) Terphenyl-d14	19.821	244	4036503	66.104	ng	0.01
Target Compounds						
						Qvalue
10) Phenol	6.893	94	89065	8.542	ng	94
15) Benzyl Alcohol	7.910	79	33205	4.325	ng	95
22) Acetophenone	8.475	105	1237431	88.772	ng	# 99
31) Naphthalene	10.475	128	188810	6.529	ng	# 97
37) 2-Methylnaphthalene	12.092	142	316675	16.912	ng	100
38) 1-Methylnaphthalene	12.316	142	288770	14.413	ng	99
46) 1,1'-Biphenyl	13.116	154	57770	2.081	ng	99
52) Acenaphthene	14.363	154	47286	2.322	ng	95
58) Fluorene	15.369	166	56268	2.125	ng	# 96
60) Diethylphthalate	15.133	149	737788	25.498	ng	99
71) Phenanthrene	17.145	178	151507	3.458	ng	98
73) Carbazole	17.527	167	144150	3.613	ng	99

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

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