

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063021\
 Data File : BP006125.D
 Acq On : 01 Jul 2021 01:04
 Operator : CG/JU
 Sample : M2896-03 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FS-02

Manual Integrations
 APPROVED

mohammad
 7/1/2021 2:53:04 PM

Quant Time: Jul 01 03:10:05 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 30 16:39:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	135172	20.000	ng	0.00
21) Naphthalene-d8	10.693	136	527974	20.000	ng	0.00
39) Acenaphthene-d10	14.528	164	321221	20.000	ng	0.00
64) Phenanthrene-d10	17.275	188	581164	20.000	ng	0.00
76) Chrysene-d12	21.351	240	469582	20.000	ng	# 0.00
86) Perylene-d12	23.757	264	531070	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	41241	4.896	ng	0.00
7) Phenol-d6	7.087	99	52397	4.799	ng	0.01
23) Nitrobenzene-d5	9.063	82	32545	3.022	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	25492	4.624	ng	0.00
45) 2-Fluorobiphenyl	13.151	172	78452	3.207	ng	0.00
79) Terphenyl-d14	19.822	244	104133	4.152	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.116	94	588389	53.944	ng	97
17) 2-Methylphenol	8.363	107	550140	74.026	ng	97
20) 3+4-Methylphenols	8.704	107	1123972	111.964	ng	97
22) Acetophenone	8.728	105	60058	4.322	ng	# 89
27) 2,4-Dimethylphenol	9.893	122	734876m	92.450	ng	
31) Naphthalene	10.746	128	223839	7.327	ng	99
32) Benzoic acid	10.051	122	13688	8.315	ng	# 82
37) 2-Methylnaphthalene	12.357	142	207549	9.543	ng	98
38) 1-Methylnaphthalene	12.575	142	122110m	5.960	ng	
58) Fluorene	15.575	166	68131	2.699	ng	# 97
71) Phenanthrene	17.322	178	80574	2.309	ng	# 92

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

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