

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011421\
 Data File : BP004557.D
 Acq On : 14 Jan 2021 15:13
 Operator : CG/JU
 Sample : M1123-01RE
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 355RE

Manual Integrations
 APPROVED

mohammad
 1/15/2021 10:08:31 AM

Quant Time: Jan 14 15:46:17 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 12 15:19:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.811	152	211243	20.00	ng	0.00
21) Naphthalene-d8	10.605	136	773984	20.00	ng	# 0.00
39) Acenaphthene-d10	14.463	164	458316	20.00	ng	0.00
64) Phenanthrene-d10	17.222	188	916193	20.00	ng	0.00
76) Chrysene-d12	21.322	240	1004058	20.00	ng	0.00
86) Perylene-d12	23.674	264	1164735	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.405	112	9449	0.78	ng	0.00
7) Phenol-d6	6.987	99	330726	19.92	ng	0.00
23) Nitrobenzene-d5	8.969	82	849283	62.04	ng	0.00
42) 2,4,6-Tribromophenol	15.969	330	1343	0.17	ng	0.00
45) 2-Fluorobiphenyl	13.081	172	2211584	62.73	ng	0.00
79) Terphenyl-d14	19.792	244	2908125	52.92	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.916	163	136834	3.60	ng	# 95
71) Phenanthrene	17.269	178	157805	3.00	ng	# 96
75) Fluoranthene	19.245	202	318048	4.98	ng	99
78) Pyrene	19.592	202	353524	5.17	ng	98
81) Benzo(a)anthracene	21.304	228	181486	2.66	ng	97
83) Chrysene	21.357	228	171787	2.64	ng	97
84) Bis(2-ethylhexyl)phtha...	21.222	149	188464	4.78	ng	99
88) Benzo(b)fluoranthene	22.963	252	248419m	3.30	ng	
90) Benzo(a)pyrene	23.574	252	173767	2.47	ng	# 92
92) Benzo(g,h,i)perylene	26.845	276	172483	2.36	ng	96

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

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