

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112021\
 Data File : BP008057.D
 Acq On : 20 Nov 2021 14:54
 Operator : CG/JU
 Sample : M4731-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C-1600-FLOOD-WALL-JG

Quant Time: Nov 22 05:31:24 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 20 00:21:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	104066	20.000	ng	0.00
21) Naphthalene-d8	10.616	136	381451	20.000	ng	-0.01
39) Acenaphthene-d10	14.463	164	240528	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	532812	20.000	ng	-0.01
76) Chrysene-d12	21.310	240	610105	20.000	ng	-0.01
86) Perylene-d12	23.704	264	709478	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.417	112	897838	147.467	ng	0.00
7) Phenol-d6	7.005	99	1078537	127.795	ng	0.00
23) Nitrobenzene-d5	8.981	82	804078	95.627	ng	-0.01
42) 2,4,6-Tribromophenol	15.957	330	444056	145.882	ng	-0.01
45) 2-Fluorobiphenyl	13.087	172	1542959	100.447	ng	0.00
79) Terphenyl-d14	19.786	244	2775372	89.437	ng	-0.01
Target Compounds						
						Qvalue
60) Diethylphthalate	15.287	149	87258	4.670	ng	98
74) Di-n-butylphthalate	18.181	149	141680	4.504	ng	99
80) Butylbenzylphthalate	20.463	149	36924	2.133	ng	98
81) Benzo(a)anthracene	21.298	228	99580	2.434	ng	100
83) Chrysene	21.345	228	108808	2.802	ng	96
84) Bis(2-ethylhexyl)phtha...	21.227	149	58345	2.492	ng	99
85) Di-n-octyl phthalate	22.151	149	131055	3.054	ng	# 85
87) Indeno(1,2,3-cd)pyrene	26.198	276	113861	2.313	ng	97
88) Benzo(b)fluoranthene	22.974	252	214218	4.829	ng	99
89) Benzo(k)fluoranthene	23.021	252	233642	5.294	ng	98
90) Benzo(a)pyrene	23.598	252	209684	5.640	ng	99

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

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