

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
 Data File : BP005268.D
 Acq On : 12 Apr 2021 16:17
 Operator : CG/JU
 Sample : M1967-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB1-C

Quant Time: Apr 12 16:57:14 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 09 18:44:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	29080	20.00	ng	# 0.00
21) Naphthalene-d8	10.463	136	98638	20.00	ng	# 0.00
39) Acenaphthene-d10	14.339	164	58349	20.00	ng	0.00
64) Phenanthrene-d10	17.104	188	143391	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	171893	20.00	ng	0.00
86) Perylene-d12	23.474	264	175587	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.281	112	163534	59.02	ng	0.00
7) Phenol-d6	6.863	99	276874	87.57	ng	0.00
23) Nitrobenzene-d5	8.840	82	192480	61.54	ng	0.00
42) 2,4,6-Tribromophenol	15.839	330	5430	7.37	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	271259	62.76	ng	0.00
79) Terphenyl-d14	19.680	244	489308	55.92	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.893	94	258573	79.86	ng	74
75) Fluoranthene	19.127	202	32532	3.27	ng	98
78) Pyrene	19.480	202	36999	3.46	ng	97
81) Benzo(a)anthracene	21.186	228	24659	2.17	ng	97
83) Chrysene	21.239	228	23778	2.16	ng	97
88) Benzo(b)fluoranthene	22.786	252	30009	2.61	ng	# 97
90) Benzo(a)pyrene	23.380	252	26622	2.50	ng	# 98

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

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