

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
 Data File : BP005269.D
 Acq On : 12 Apr 2021 16:51
 Operator : CG/JU
 Sample : M1967-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB2-B

Manual Integrations
 APPROVED

mohammad
 4/14/2021 11:49:23 AM

Quant Time: Apr 12 19:18:42 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	28052	20.00	ng	# 0.00
21) Naphthalene-d8	10.463	136	95051	20.00	ng	# 0.00
39) Acenaphthene-d10	14.340	164	57155	20.00	ng	0.00
64) Phenanthrene-d10	17.104	188	135090	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	168026	20.00	ng	0.00
86) Perylene-d12	23.474	264	174828	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.275	112	59902	22.41	ng	0.00
7) Phenol-d6	6.864	99	207234	67.95	ng	0.00
23) Nitrobenzene-d5	8.834	82	164708	54.65	ng	0.00
42) 2,4,6-Tribromophenol	15.840	330	534	0.74	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	210905	49.82	ng	0.00
79) Terphenyl-d14	19.681	244	382958	44.77	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.887	94	83992	26.89	ng	72
50) Dimethylphthalate	13.798	163	9688	2.31	ng	# 98
75) Fluoranthene	19.128	202	38299	4.09	ng	96
78) Pyrene	19.481	202	41699	3.99	ng	99
81) Benzo(a)anthracene	21.186	228	26122	2.36	ng	97
83) Chrysene	21.239	228	27787	2.58	ng	96
84) Bis(2-ethylhexyl)phtha...	21.116	149	17000	2.77	ng	95
88) Benzo(b)fluoranthene	22.786	252	37028m	3.24	ng	
90) Benzo(a)pyrene	23.374	252	28818	2.72	ng	# 90

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

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