

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041621\
 Data File : BP005321.D
 Acq On : 16 Apr 2021 16:32
 Operator : CG/JU
 Sample : M1947-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FS-SB19B(19-20)X

Quant Time: Apr 16 17:38:25 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 16 11:02:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.658	152	30159	20.00	ng	0.00
21) Naphthalene-d8	10.452	136	74125	20.00	ng	0.00
39) Acenaphthene-d10	14.328	164	47864	20.00	ng	0.00
64) Phenanthrene-d10	17.092	188	115621	20.00	ng	0.00
76) Chrysene-d12	21.192	240	179160	20.00	ng	0.00
86) Perylene-d12	23.463	264	179492	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.264	112	260318	105.25	ng	0.00
7) Phenol-d6	6.846	99	292636	102.30	ng	0.00
23) Nitrobenzene-d5	8.822	82	186726	77.80	ng	0.00
42) 2,4,6-Tribromophenol	15.834	330	78564	123.48	ng	0.00
45) 2-Fluorobiphenyl	12.940	172	302999	84.08	ng	0.00
79) Terphenyl-d14	19.669	244	648108	71.45	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.116	202	22708	2.79	ng	94
78) Pyrene	19.469	202	24735	2.42	ng	96
81) Benzo(a)anthracene	21.180	228	26828	2.35	ng	99
83) Chrysene	21.228	228	22641	2.00	ng	98
88) Benzo(b)fluoranthene	22.775	252	34073	2.83	ng	98
90) Benzo(a)pyrene	23.363	252	25868	2.35	ng	# 99

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

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