

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092721\
 Data File : BP007216.D
 Acq On : 27 Sep 2021 21:14
 Operator : CG/JU
 Sample : M3950-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HD-02-092421

Quant Time: Sep 28 02:51:57 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 23 11:17:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	220524	20.000	ng	-0.01
21) Naphthalene-d8	10.681	136	929536	20.000	ng	#-0.01
39) Acenaphthene-d10	14.516	164	672590	20.000	ng	-0.01
64) Phenanthrene-d10	17.269	188	1581604	20.000	ng	0.00
76) Chrysene-d12	21.351	240	1284743	20.000	ng	0.00
86) Perylene-d12	23.768	264	1258538	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.458	112	1714894	130.173	ng	0.00
7) Phenol-d6	7.063	99	2175587	120.829	ng	0.00
23) Nitrobenzene-d5	9.046	82	1264381	79.215	ng	-0.01
42) 2,4,6-Tribromophenol	16.010	330	1445704	165.796	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	3593908	76.551	ng	0.00
79) Terphenyl-d14	19.827	244	6306512	94.065	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.275	202	321083	3.286	ng	97
78) Pyrene	19.627	202	308764	3.663	ng	99
83) Chrysene	21.386	228	160178	2.010	ng	# 91
88) Benzo(b)fluoranthene	23.033	252	224882	2.990	ng	# 83
90) Benzo(a)pyrene	23.657	252	168891	2.643	ng	# 85

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

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