

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070823\
 Data File : BP016136.D
 Acq On : 09 Jul 2023 17:22
 Operator : MA/JU
 Sample : PB153808BL
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK808

Quant Time: Jul 09 22:43:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 08 10:11:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	190907	20.000	ng/ul	0.00
20) Naphthalene-d8	10.851	136	822001	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.675	164	515832	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	1059999	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.563	240	835687	20.000	ng/ul	# 0.02
88) Perylene-d12	24.092	264	779337	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	35412	6.674	ng/uL	0.00
4) Pyridine-d5	3.799	84	408418	30.534	ng/ul	0.00
7) Phenol-d5	7.228	99	468624	28.690	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.352	67	267541	26.474	ng/ul	0.00
11) 2-Chlorophenol-d4	7.563	132	407138	33.019	ng/ul	0.00
15) 4-Methylphenol-d8	8.781	113	363901	28.201	ng/ul	0.01
21) Nitrobenzene-d5	9.228	128	196915	31.346	ng/ul	0.01
24) 2-Nitrophenol-d4	9.975	143	210988	28.777	ng/ul	0.03
28) 2,4-Dichlorophenol-d3	10.563	165	263681	20.182	ng/ul	0.06
31) 4-Chloroaniline-d4	11.063	131	459304	27.366	ng/ul	0.04
46) Dimethylphthalate-d6	14.098	166	1376764	34.531	ng/ul	0.01
49) Acenaphthylene-d8	14.369	160	1465289	32.693	ng/ul	0.00
54) 4-Nitrophenol-d4	15.057	143	22158	4.211	ng/ul	0.09
60) Fluorene-d10	15.686	176	1113230	33.910	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.886	200	109425	16.786	ng/ul	0.04
73) Anthracene-d10	17.551	188	1671155	34.515	ng/ul	0.00
81) Pyrene-d10	19.780	212	1942439	35.596	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.915	264	1366900	34.769	ng/ul	0.00

Target Compounds	Qvalue

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

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