

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
 Data File : BP018733.D
 Acq On : 11 Dec 2023 18:51
 Operator : MA/JU
 Sample : PB157686BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK686

Quant Time: Dec 11 22:01:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 11 22:02:13 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	298348	20.000	ng/ul	0.00
20) Naphthalene-d8	10.634	136	931310	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	716105	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	1645621	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	1696858	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	1952766	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	45595	5.208	ng/uL	0.00
4) Pyridine-d5	3.681	84	605152	25.805	ng/ul	0.00
7) Phenol-d5	7.011	99	727594	24.301	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	435895	24.564	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	577086	24.702	ng/ul	0.00
15) 4-Methylphenol-d8	8.552	113	567350	23.415	ng/ul	0.00
21) Nitrobenzene-d5	9.005	128	266583	32.161	ng/ul	0.00
24) 2-Nitrophenol-d4	9.722	143	230456	26.817	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	452666	25.760	ng/ul	0.00
31) 4-Chloroaniline-d4	10.775	131	627532	27.326	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	1724309	26.979	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	1890083	26.857	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	269536	25.692	ng/ul	0.00
60) Fluorene-d10	15.463	176	1480881	27.603	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	215439	21.834	ng/ul	0.00
73) Anthracene-d10	17.316	188	2391916	27.648	ng/ul	0.00
81) Pyrene-d10	19.557	212	2967829	22.469	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.521	264	3112657	27.338	ng/ul	0.00

Target Compounds Qvalue

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

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