

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018316.D
 Acq On : 23 Nov 2023 14:59
 Operator : MA/JU
 Sample : PB157189BL
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK189

Quant Time: Nov 24 00:10:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 22 14:09:55 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	451832	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	1868142	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.504	164	1173522	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	2155005	20.000	ng/ul	-0.01
79) Chrysene-d12	21.357	240	1402779	20.000	ng/ul	0.00
88) Perylene-d12	23.780	264	1752270	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	77612	6.434	ng/uL	0.00
4) Pyridine-d5	3.687	84	1034217	31.196	ng/ul	0.00
7) Phenol-d5	7.034	99	1269570	30.745	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	766058	31.024	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.399	132	1058153	30.521	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	1035337	31.069	ng/ul	-0.01
21) Nitrobenzene-d5	9.034	128	514609	33.915	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	502949	35.890	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.287	165	1029138	30.440	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.810	131	1542012	33.817	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	3270334	31.319	ng/ul	-0.01
49) Acenaphthylene-d8	14.198	160	3753184	31.544	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	376739	26.135	ng/ul	-0.01
60) Fluorene-d10	15.498	176	2689737	30.796	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	240615	26.229	ng/ul	-0.01
73) Anthracene-d10	17.351	188	3669464	31.857	ng/ul	-0.01
81) Pyrene-d10	19.592	212	3388540	30.351	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.621	264	3264519	32.135	ng/ul	0.00

Target Compounds Qvalue

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

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