

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010524\
 Data File : BP019279.D
 Acq On : 05 Jan 2024 12:26
 Operator : MA/JU
 Sample : PB158131BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK131

Quant Time: Jan 05 23:55:15 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	130124	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.592	136	494731	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	329364	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	804951	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	941118	20.000	ng/ul	0.00
88) Perylene-d12	23.668	264	1119745	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	19666	5.162	ng/uL	0.00
4) Pyridine-d5	3.652	84	273037	28.551	ng/ul	0.00
7) Phenol-d5	6.987	99	312987	26.822	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	182347	27.011	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	245928	26.142	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	242866	26.733	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	119177	27.867	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	137693	29.791	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	242112	25.941	ng/ul	0.00
31) 4-Chloroaniline-d4	10.745	131	330467	28.786	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	808833	27.803	ng/ul	-0.01
49) Acenaphthylene-d8	14.133	160	880298	27.887	ng/ul	0.00
54) 4-Nitrophenol-d4	14.692	143	138058	28.103	ng/ul	0.03
60) Fluorene-d10	15.439	176	697507	27.753	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	121081	22.945	ng/ul	0.00
73) Anthracene-d10	17.298	188	1199805	28.542	ng/ul	0.00
81) Pyrene-d10	19.545	212	1565015	27.286	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	1793751	27.869	ng/ul	0.00

Target Compounds Qvalue

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

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