

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032524\
 Data File : BP019978.D
 Acq On : 25 Mar 2024 10:14
 Operator : MA/JU
 Sample : PB159425BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK425

Quant Time: Mar 25 10:59:15 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 10:46:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	108609	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.346	136	432697	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.239	164	279020	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.063	188	588862	20.000	ng/u1	-0.01
79) Chrysene-d12	21.539	240	511849	20.000	ng/u1	0.00
88) Perylene-d12	24.827	264	564326	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	17788	7.176	ng/uL	-0.02
4) Pyridine-d5	3.622	84	249130	33.285	ng/u1	-0.02
7) Phenol-d5	6.787	99	306780	33.491	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.946	67	202752	33.974	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.134	132	219893	33.427	ng/u1	-0.02
15) 4-Methylphenol-d8	8.304	113	234829	33.136	ng/u1	0.00
21) Nitrobenzene-d5	8.746	128	112770	34.002	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.451	143	128795	34.709	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	9.987	165	199777	29.479	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.504	131	305575	34.589	ng/u1	-0.01
46) Dimethylphthalate-d6	13.651	166	680860	34.804	ng/u1	0.00
49) Acenaphthylene-d8	13.922	160	787847	34.475	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.492	143	87806	28.551	ng/u1	0.01
60) Fluorene-d10	15.263	176	574088	34.450	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	98056	27.818	ng/u1	-0.02
73) Anthracene-d10	17.169	188	882205	33.844	ng/u1	0.00
81) Pyrene-d10	19.563	212	1023796	34.091	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.592	264	947653	34.954	ng/u1	-0.02

Target Compounds Qvalue

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

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