

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102323\
 Data File : BP017780.D
 Acq On : 24 Oct 2023 16:34
 Operator : MA/JU
 Sample : PB156501BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK501

Quant Time: Oct 25 03:25:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	127104	20.000	ng/u1	0.00
20) Naphthalene-d8	10.716	136	537661	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	323165	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	675231	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.516	240	530090	20.000	ng/u1	0.00
88) Perylene-d12	24.068	264	544099	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	21943	6.422	ng/uL	0.00
4) Pyridine-d5	3.687	84	249077	27.844	ng/u1	0.00
7) Phenol-d5	7.040	99	325813	30.804	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	215925	33.056	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	266822	30.930	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	272499	32.807	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	135528	31.108	ng/u1	0.00
24) 2-Nitrophenol-d4	9.804	143	145886	30.306	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	255784	27.088	ng/u1	0.00
31) 4-Chloroaniline-d4	10.893	131	324723	26.277	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	816984	30.115	ng/u1	0.00
49) Acenaphthylene-d8	14.275	160	912281	30.154	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	85308	22.069	ng/u1	0.00
60) Fluorene-d10	15.586	176	689992	29.423	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	92904	20.888	ng/u1	0.00
73) Anthracene-d10	17.475	188	1032285	29.916	ng/u1	0.00
81) Pyrene-d10	19.733	212	1142459	34.376	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.898	264	917440	30.274	ng/u1	0.00

Target Compounds Qvalue

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

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