

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007457.D
 Acq On : 18 Oct 2021 11:14
 Operator : CG/JU
 Sample : PB139922BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK922

Quant Time: Oct 18 11:43:54 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 14 13:04:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	183708	20.00	ng/ul	-0.01
20) Naphthalene-d8	10.763	136	737278	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.598	164	477436	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	1040395	20.00	ng/ul	0.00
79) Chrysene-d12	21.439	240	1084083	20.00	ng/ul	0.00
88) Perylene-d12	23.898	264	1125247	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	29905	6.43	ng/uL	0.00
4) Pyridine-d5	3.793	84	399298	31.36	ng/ul	0.00
7) Phenol-d5	7.116	99	479305	33.41	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	280449	32.77	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	380131	33.68	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	378202	34.22	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	181073	38.81	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	184071	42.24	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.381	165	359670	33.84	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	508730	35.39	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	1210335	36.64	ng/ul	-0.01
49) Acenaphthylene-d8	14.287	160	1434216	34.55	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	196417	37.06	ng/ul	0.00
60) Fluorene-d10	15.592	176	1004546	36.34	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	135355	32.47	ng/ul	0.00
73) Anthracene-d10	17.451	188	1678994	38.00	ng/ul	0.00
81) Pyrene-d10	19.681	212	1996922	36.85	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.739	264	1984726	35.45	ng/ul	0.00

Target Compounds	Qvalue

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

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