

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091521\
 Data File : BP006978.D
 Acq On : 15 Sep 2021 16:47
 Operator : CG/JU
 Sample : PB138901BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK901

Quant Time: Sep 15 17:16:01 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 15 17:01:19 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	424322	20.000	ng/u1	0.00
20) Naphthalene-d8	10.722	136	1745690	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.545	164	1132238	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.298	188	2272019	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.374	240	1849662	20.000	ng/u1	# 0.00
88) Perylene-d12	23.798	264	1635068	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	68188	6.395	ng/uL	0.00
4) Pyridine-d5	3.770	84	855749	30.798	ng/u1	0.00
7) Phenol-d5	7.081	99	1019859	32.702	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	608992	34.001	ng/u1	0.00
11) 2-Chlorophenol-d4	7.452	132	844559	33.545	ng/u1	0.00
15) 4-Methylphenol-d8	8.622	113	816084	33.323	ng/u1	0.00
21) Nitrobenzene-d5	9.081	128	393645	37.228	ng/u1	0.00
24) 2-Nitrophenol-d4	9.805	143	409264	39.314	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	788862	32.138	ng/u1	0.00
31) 4-Chloroaniline-d4	10.857	131	1154494	34.011	ng/u1	0.00
46) Dimethylphthalate-d6	13.957	166	2580079	35.183	ng/u1	0.00
49) Acenaphthylene-d8	14.240	160	3141775	33.827	ng/u1	0.00
54) 4-Nitrophenol-d4	14.734	143	377216	29.609	ng/u1	0.00
60) Fluorene-d10	15.540	176	2224181	34.366	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	301428	29.990	ng/u1	0.00
73) Anthracene-d10	17.392	188	3378364	35.916	ng/u1	0.00
81) Pyrene-d10	19.628	212	3609026	41.663	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.639	264	2728973	34.597	ng/u1	0.00

Target Compounds Qvalue

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

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