

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006943.D
 Acq On : 10 Sep 2021 16:19
 Operator : CG/JU
 Sample : PB138947BL
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK947

Quant Time: Sep 10 17:06:46 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 09 11:07:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	354869	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	1419527	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.563	164	879289	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.310	188	1782425	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.392	240	1731717	20.000	ng/u1	# 0.00
88) Perylene-d12	23.827	264	1823507	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	51937	5.824	ng/uL	0.00
4) Pyridine-d5	3.787	84	686653	29.549	ng/u1	0.00
7) Phenol-d5	7.093	99	786070	30.138	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	461615	30.817	ng/u1	0.00
11) 2-Chlorophenol-d4	7.464	132	641932	30.487	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	611415	29.852	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	285316	33.183	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	280949	33.189	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	576598	28.887	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	838061	30.361	ng/u1	0.00
46) Dimethylphthalate-d6	13.969	166	1811196	31.803	ng/u1	0.00
49) Acenaphthylene-d8	14.251	160	2195538	30.440	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	267980	27.086	ng/u1	0.00
60) Fluorene-d10	15.551	176	1566299	31.163	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	177470	22.507	ng/u1	0.00
73) Anthracene-d10	17.404	188	2452687	33.237	ng/u1	-0.01
81) Pyrene-d10	19.639	212	2837783	34.991	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.669	264	2745652	31.211	ng/u1	-0.01

Target Compounds Qvalue

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

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