

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006921.D
 Acq On : 10 Sep 2021 03:47
 Operator : CG/JU
 Sample : PB138947BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK947

Quant Time: Sep 10 04:26:06 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	357833	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	1425514	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.563	164	882006	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.310	188	1823831	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.386	240	1783265	20.000	ng/u1	-0.01
88) Perylene-d12	23.821	264	1840771	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	55927	6.220	ng/uL	0.00
4) Pyridine-d5	3.787	84	738607	31.521	ng/u1	0.00
7) Phenol-d5	7.093	99	876198	33.316	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	525171	34.769	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	706407	33.271	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	671086	32.494	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	310318	35.939	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	306255	36.027	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	622627	31.063	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	926946	33.441	ng/u1	0.00
46) Dimethylphthalate-d6	13.969	166	1993892	34.904	ng/u1	0.00
49) Acenaphthylene-d8	14.251	160	2402511	33.207	ng/u1	0.00
54) 4-Nitrophenol-d4	14.739	143	302004	30.431	ng/u1	0.00
60) Fluorene-d10	15.551	176	1707340	33.864	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	216363	26.817	ng/u1	0.00
73) Anthracene-d10	17.404	188	2774488	36.745	ng/u1	-0.01
81) Pyrene-d10	19.639	212	3211630	38.456	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.668	264	3129217	35.237	ng/u1	-0.01

Target Compounds	Qvalue

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

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