

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091321\
 Data File : BP006964.D
 Acq On : 13 Sep 2021 17:24
 Operator : CG/JU
 Sample : PB139008BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK008

Quant Time: Sep 13 18:04:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 13 14:04:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	419745	20.000	ng/u1	0.00
20) Naphthalene-d8	10.722	136	1776642	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.551	164	1205030	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.304	188	2534042	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.386	240	2334353	20.000	ng/u1	# 0.00
88) Perylene-d12	23.815	264	2309055	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	75766	7.183	ng/uL	0.00
4) Pyridine-d5	3.764	84	910203	33.115	ng/u1	0.00
7) Phenol-d5	7.075	99	1112920	36.075	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	659534	37.224	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.446	132	921335	36.994	ng/u1	0.00
15) 4-Methylphenol-d8	8.622	113	898448	37.086	ng/u1	0.00
21) Nitrobenzene-d5	9.081	128	442488	41.118	ng/u1	0.00
24) 2-Nitrophenol-d4	9.799	143	469452	44.310	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	883014	35.347	ng/u1	0.00
31) 4-Chloroaniline-d4	10.857	131	1292594	37.416	ng/u1	0.00
46) Dimethylphthalate-d6	13.963	166	3047865	39.052	ng/u1	0.00
49) Acenaphthylene-d8	14.245	160	3623621	36.659	ng/u1	0.00
54) 4-Nitrophenol-d4	14.734	143	486846	35.906	ng/u1	0.00
60) Fluorene-d10	15.545	176	2587510	37.564	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.657	200	403566	36.001	ng/u1	0.00
73) Anthracene-d10	17.404	188	3971355	37.855	ng/u1	0.00
81) Pyrene-d10	19.633	212	4415456	40.389	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.657	264	4195361	37.662	ng/u1	0.00

Target Compounds Qvalue

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

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